

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1623kxg

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

| | | | |
|------|----|--------|--|
| NEWS | 1 | | Web Page URLs for STN Seminar Schedule - N. America |
| NEWS | 2 | Apr 08 | "Ask CAS" for self-help around the clock |
| NEWS | 3 | Apr 09 | BEILSTEIN: Reload and Implementation of a New Subject Area |
| NEWS | 4 | Apr 09 | ZDB will be removed from STN |
| NEWS | 5 | Apr 19 | US Patent Applications available in IFICDB, IFIPAT, and IFIUDB |
| NEWS | 6 | Apr 22 | Records from IP.com available in CAPLUS, HCAPLUS, and ZCAPLUS |
| NEWS | 7 | Apr 22 | BIOSIS Gene Names now available in TOXCENTER |
| NEWS | 8 | Apr 22 | Federal Research in Progress (FEDRIP) now available |
| NEWS | 9 | Jun 03 | New e-mail delivery for search results now available |
| NEWS | 10 | Jun 10 | MEDLINE Reload |
| NEWS | 11 | Jun 10 | PCTFULL has been reloaded |
| NEWS | 12 | Jul 02 | FOREGE no longer contains STANDARDS file segment |
| NEWS | 13 | Jul 22 | USAN to be reloaded July 28, 2002; saved answer sets no longer valid |
| NEWS | 14 | Jul 29 | Enhanced polymer searching in REGISTRY |
| NEWS | 15 | Jul 30 | NETFIRST to be removed from STN |
| NEWS | 16 | Aug 08 | CANCERLIT reload |
| NEWS | 17 | Aug 08 | PHARMAMarketLetter(PHARMAML) - new on STN |
| NEWS | 18 | Aug 08 | NTIS has been reloaded and enhanced |
| NEWS | 19 | Aug 19 | Aquatic Toxicity Information Retrieval (AQUIRE) now available on STN |
| NEWS | 20 | Aug 19 | IFIPAT, IFICDB, and IFIUDB have been reloaded |
| NEWS | 21 | Aug 19 | The MEDLINE file segment of TOXCENTER has been reloaded |
| NEWS | 22 | Aug 26 | Sequence searching in REGISTRY enhanced |
| NEWS | 23 | Sep 03 | JAPIO has been reloaded and enhanced |
| NEWS | 24 | Sep 16 | Experimental properties added to the REGISTRY file |
| NEWS | 25 | Sep 16 | CA Section Thesaurus available in CAPLUS and CA |
| NEWS | 26 | Oct 01 | CASREACT Enriched with Reactions from 1907 to 1985 |
| NEWS | 27 | Oct 21 | EVENTLINE has been reloaded |
| NEWS | 28 | Oct 24 | BEILSTEIN adds new search fields |
| NEWS | 29 | Oct 24 | Nutraceuticals International (NUTRACEUT) now available on STN |
| NEWS | 30 | Oct 25 | MEDLINE SDI run of October 8, 2002 |
| NEWS | 31 | Nov 18 | DKILIT has been renamed APOLLIT |
| NEWS | 32 | Nov 25 | More calculated properties added to REGISTRY |
| NEWS | 33 | Dec 02 | TIBKAT will be removed from STN |
| NEWS | 34 | Dec 04 | CSA files on STN |
| NEWS | 35 | Dec 17 | PCTFULL now covers WP/PCT Applications from 1978 to date |
| NEWS | 36 | Dec 17 | TOXCENTER enhanced with additional content |
| NEWS | 37 | Dec 17 | Adis Clinical Trials Insight now available on STN |
| NEWS | 38 | Dec 30 | ISMEC no longer available |
| NEWS | 39 | Jan 13 | Indexing added to some pre-1967 records in CA/CAPLUS |
| NEWS | 40 | Jan 21 | NUTRACEUT offering one free connect hour in February 2003 |
| NEWS | 41 | Jan 21 | PHARMAML offering one free connect hour in February 2003 |
| NEWS | 42 | Jan 29 | Simultaneous left and right truncation added to COMPENDEX, ENERGY, INSPEC |
| NEWS | 43 | Feb 13 | CANCERLIT is no longer being updated |
| NEWS | 44 | Feb 24 | METADEX enhancements |
| NEWS | 45 | Feb 24 | PCTGEN now available on STN |

NEWS 46 Feb 24 TEMA now available on STN
NEWS 47 Feb 26 NTIS now allows simultaneous left and right truncation
NEWS 48 Feb 26 PCTFULL now contains images
NEWS 49 Mar 04 SDI PACKAGE for monthly delivery of multifile SDI results

NEWS EXPRESS January 6 CURRENT WINDOWS VERSION IS V6.01a,
CURRENT MACINTOSH VERSION IS V6.0b(ENG) AND V6.0Jb(JP),
AND CURRENT DISCOVER FILE IS DATED 01 OCTOBER 2002
NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS INTER General Internet Information
NEWS LOGIN Welcome Banner and News Items
NEWS PHONE Direct Dial and Telecommunication Network Access to STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 15:02:14 ON 17 MAR 2003

=> le reg

LE IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system.
For a list of commands available to you in the current file, enter "HELP COMMANDS" at an arrow prompt (=>).

=> file reg

| COST IN U.S. DOLLARS | SINCE FILE ENTRY | TOTAL SESSION |
|----------------------|------------------|---------------|
| FULL ESTIMATED COST | 0.21 | 0.21 |

FILE 'REGISTRY' ENTERED AT 15:02:24 ON 17 MAR 2003

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2003 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 16 MAR 2003 HIGHEST RN 499182-00-2

DICTIONARY FILE UPDATES: 16 MAR 2003 HIGHEST RN 499182-00-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

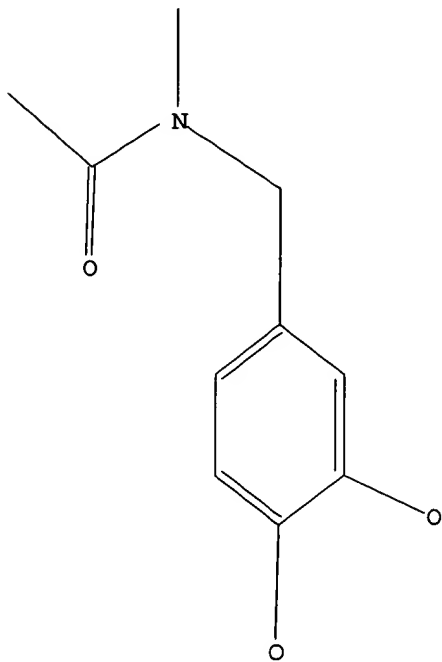
Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:

<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=>
Uploading 09787764-1.str

L1 STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1 exact sam
SAMPLE SEARCH INITIATED 15:04:04 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA EXA SAM L1

=> s l1 sss sam
SAMPLE SEARCH INITIATED 15:04:49 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 373 TO ITERATE

100.0% PROCESSED 373 ITERATIONS 4 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**
PROJECTED ITERATIONS: 6302 TO 8618
PROJECTED ANSWERS: 4 TO 200

4 SEA SSS SAM L1

```

L3      4 ANSWERS      REGISTRY  COPYRIGHT 2003 ACS
IN      Benzamide, 3-[2-(dimethylamino)ethoxy]-4-methoxy-N-[(1S)-3-methyl-1-
        [[[1S)-1-methyl-2-oxo-3-[(2-pyridinylsulfonyl)amino]propyl]amino]carbonyl
        ]butyl]- (9CI)
MF      C27 H39 N5 O7 S

```

CN(C)CCOC1=CC=C(C=C1)C(=O)N[C@@H](C)C(=O)N[C@H](C)C(=O)NS(=O)(=O)c2ccncc2

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1) :1

```
L3      4 ANSWERS      REGISTRY  COPYRIGHT 2003 ACS
IN      Benzeneacetamide, N- [(3S)-3- [[ (2S)-2- [[3- [2- (dimethylamino)ethoxy]-4-
        methoxybenzoyl] amino]-4-methyl-1-oxopentyl] amino]-2-oxobutyl]-.alpha.-
        methyl-3- (2-pyridinyl) - (9CI)
MF      C36 H47 N5 O6
```

CC(C(=O)NCC(=O)S[C@H](C)NC(=O)S[C@@H](C)NC(=O)c1ccc(OC)c(OCCN)c1)c2ccc(cc2)c3ccccc3n

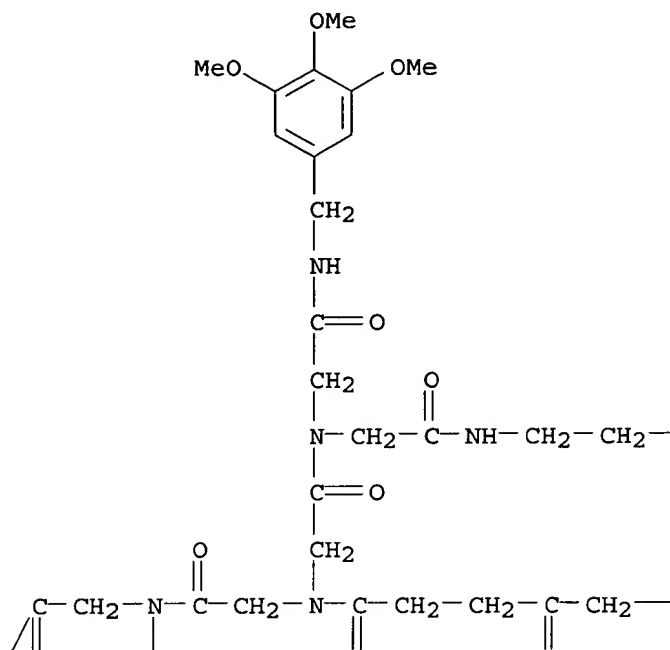
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 4 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Glycinamide, 1,1'-(1,4,7-trioxo-1,7-heptanediyl)bis[N-[2-[[2-[(4-methoxyphenyl)ethyl]amino]-2-oxoethyl][2-oxo-2-[[3,4,5-trimethoxyphenyl)methyl]amino]ethyl]amino]-2-oxoethyl]glycyl-N2-[2-[[2-(4-

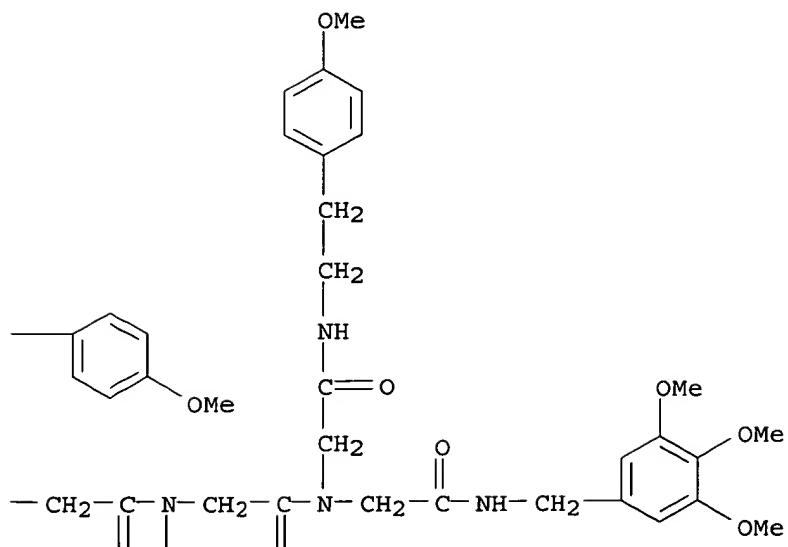
methoxyphenyl) ethyl] amino] -2-oxoethyl] -N- [(3,4,5-trimethoxyphenyl) methyl] -
(9CI)

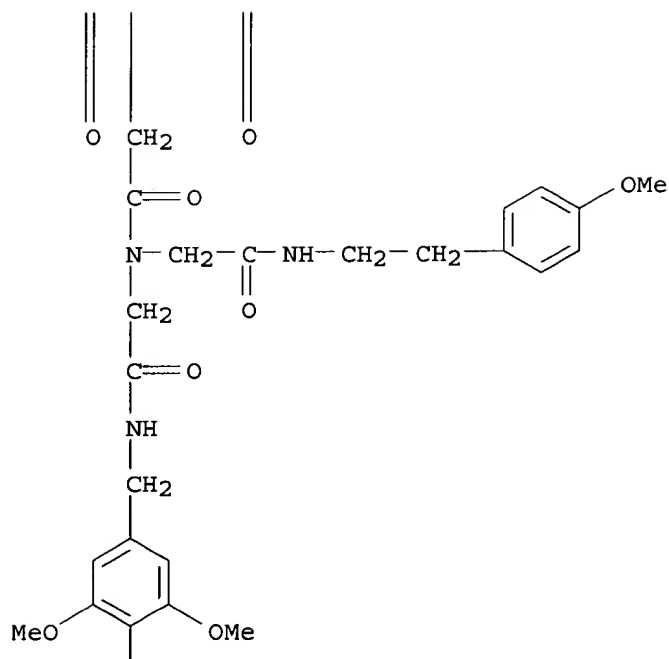
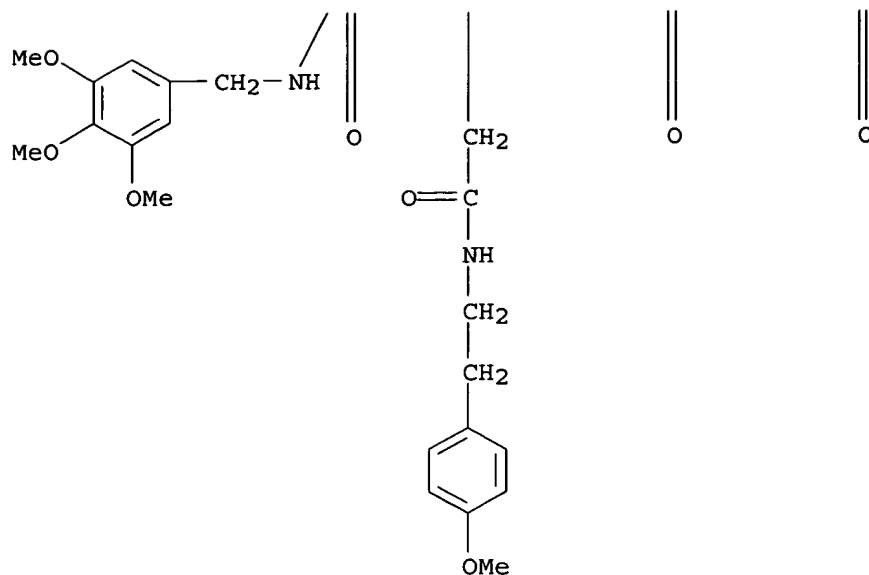
MF C107 H136 N14 O31

PAGE 1-A



PAGE 1-B



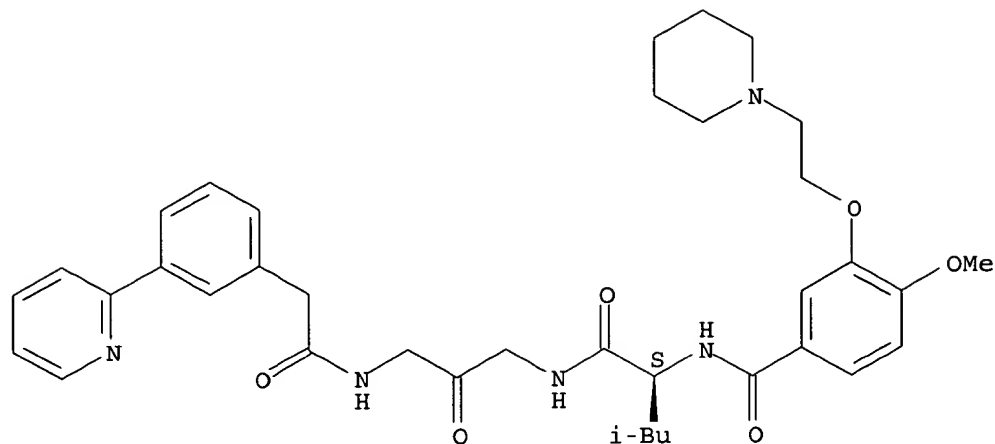


HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 4 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzeneacetamide, N-[3-[(2S)-2-[[4-methoxy-3-[2-(1-

piperidinyl)ethoxy]benzoyl]amino]-4-methyl-1-oxopentyl]amino]-2-oxopropyl]-
 3-(2-pyridinyl)-(9CI)
 MF C37 H47 N5 O6

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> s l1 sss full
 FULL SEARCH INITIATED 15:06:32 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 7741 TO ITERATE

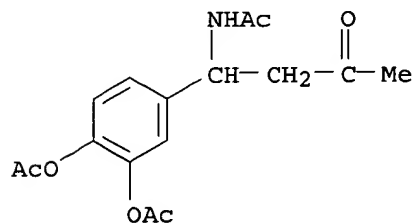
100.0% PROCESSED 7741 ITERATIONS
 SEARCH TIME: 00.00.01

81 ANSWERS

L4 81 SEA SSS FUL L1

=> d scan

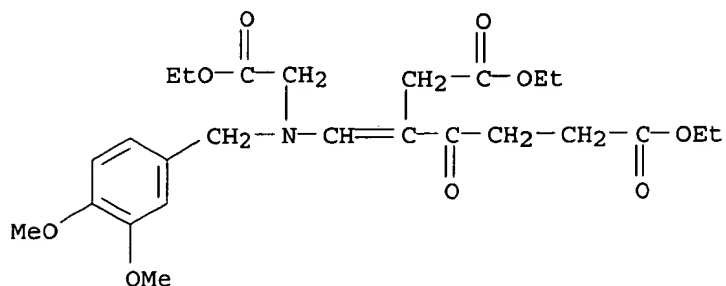
L4 81 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Acetamide, N-[1-[[3,4-bis(acetyloxy)phenyl]methyl]-2-oxopropyl]-(9CI)
 MF C16 H19 N O6



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

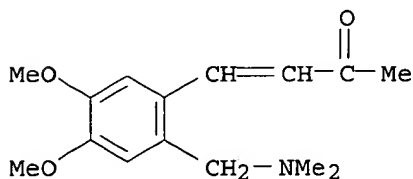
L4 81 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Heptanedioic acid, 3-[[[(3,4-dimethoxyphenyl)methyl] (2-ethoxy-2-oxoethyl)amino]methylene]-4-oxo-, diethyl ester (9CI)
MF C25 H35 N O9



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

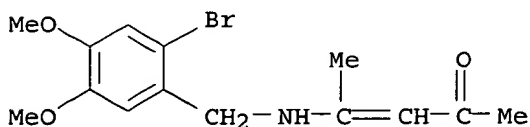
L4 81 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 3-Buten-2-one, 4-[2-[(dimethylamino)methyl]-4,5-dimethoxyphenyl]- (9CI)
MF C15 H21 N O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

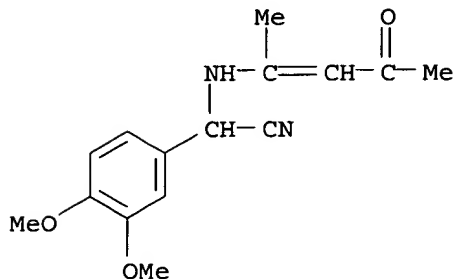
L4 81 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 3-Penten-2-one, 4-[[[2-bromo-4,5-dimethoxyphenyl)methyl]amino]- (9CI)
MF C14 H18 Br N O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L4 81 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Benzeneacetone nitrile, 3,4-dimethoxy-.alpha.-[(1-methyl-3-oxo-1-butenyl)amino]- (9CI)
 MF C15 H18 N2 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

151.75

151.96

FILE 'CAPLUS' ENTERED AT 15:08:26 ON 17 MAR 2003

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 17 Mar 2003 VOL 138 ISS 12

FILE LAST UPDATED: 16 Mar 2003 (20030316/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l4 and vannillinamide

45 L4

0 VANNILLINAMIDE

L5

0 L4 AND VANNILLINAMIDE

=> s l5 and N-acyl

2546853 N

90371 ACYL

229 ACYLS

90469 ACYL

(ACYL OR ACYLS)

7921 N-ACYL
(N(W)ACYL)

L6 0 L5 AND N-ACYL

=> s 15 and vanil?
18727 VANIL?

L7 0 L5 AND VANIL?

=> s 17 and vanillin?
11056 VANILLIN?

L8 0 L7 AND VANILLIN?

=> s 15 and vanillinamide
0 VANILLINAMIDE

L9 0 L5 AND VANILLINAMIDE

=> s 14 and vanil?
45 L4
18727 VANIL?

L10 2 L4 AND VANIL?

=> dis l10 1-2 bib abs hitstr

L10 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2003 ACS

AN 2002:511742 CAPLUS

DN 137:216814

TI N-Acylvanillamides: Development of an Expeditious Synthesis and Discovery of New Acyl Templates for Powerful Activation of the **Vanilloid** Receptor

AU Appendino, Giovanni; Minassi, Alberto; Morello, Aniello Schiano; De Petrocellis, Luciano; Di Marzo, Vincenzo

CS Dipartimento di Scienze Chimiche Alimentari, Farmaceutiche e Farmacologiche, Novara, 28100, Italy

SO Journal of Medicinal Chemistry (2002), 45(17), 3739-3745
CODEN: JMCMAR; ISSN: 0022-2623

PB American Chemical Society

DT Journal

LA English

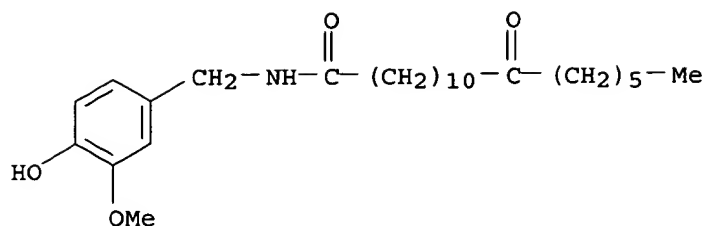
OS CASREACT 137:216814

AB A simple and general synthesis of **vanillamides** was developed and employed to screen acids from the fatty and isoprenoid pools for new acyl templates of biol. relevance as capsaicin analogs. Potent activation of the human **vanilloid** receptor 1 (VR1) was obsd. for the **vanillamides** of certain polyfunctional acids from both pools, showing that the **vanilloid** activity of capsaicinoids can be substantially improved by introducing polar groups and/or unsaturations on the acyl moiety. The activity of the unsatd. analogs was maintained or even increased by cyclopropanation, while .omega. dimerization led to a substantial increase of activity. Because of the wide structural diversity of the library of compds. screened, these observations could not be translated into a single framework of structure-activity relationships. Nevertheless, a series of new highly active leads was identified, validating the pharmacol. potential of the unnatural combination of natural building blocks to provide new bioactive compds.

IT 457067-08-2P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)
(prepn. of N-acylvanillamines as templates for **vanilloid** receptor activators)

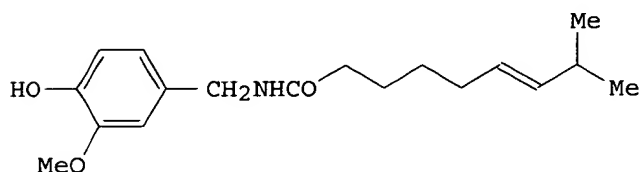
RN 457067-08-2 CAPLUS

CN Octadecanamide, N-[(4-hydroxy-3-methoxyphenyl)methyl]-12-oxo- (9CI) (CA INDEX NAME)



RE.CNT 35 THERE ARE 35 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2003 ACS
AN 1993:650209 CAPLUS
DN 119:250209
TI **Vanilloids**. 1. Analogs of capsaicin with antinociceptive and antiinflammatory activity
AU Janusz, John M.; Buckwalter, Brian L.; Young, Patricia A.; LaHann, Thomas R.; Farmer, Ralph W.; Kasting, Gerald B.; Loomans, Maurice E.; Kerckaert, Gary A.; Maddin, Cherie S.; et al.
CS Miami Valley Lab., Procter and Gamble Co., Cincinnati, OH, 45239-8707, USA
SO Journal of Medicinal Chemistry (1993), 36(18), 2595-604
CODEN: JMCMAR; ISSN: 0022-2623
DT Journal
LA English
GI



I

AB As part of a program to establish structure-activity relationships for **vanilloids**, analogs of the pungent principle capsaicin (I), the alkyl chain portion of the parent structure (and related compds. derived from homovanillic acid) was varied. In antinociceptive and antiinflammatory assays (rat and mouse hot plate and croton oil-inflamed mouse ear), compds. with widely varying alkyl chain structures were active. Short-chain compds. were active by systemic administration in the assay mentioned above but they retained the high pungency and acute toxicity characteristic of capsaicin. In contrast, the long chain cis-unsaturates, NE-19550 (**vanillyloleamide**) and NE-28345 (oleylhomovanillamide), were orally active, less pungent, and less acutely toxic than capsaicin. The potential of these compds. as antiinflammatory/analgesic agents is discussed in light of recent data on the mechanism of action of **vanilloids** on sensory nerve fibers.
IT **150988-84-4P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. and antinociceptive and antiinflammatory activity of)
RN 150988-84-4 CAPLUS
CN Octadecanamide, N-[(4-hydroxy-3-methoxyphenyl)methyl]-9-oxo- (9CI) (CA INDEX NAME)

